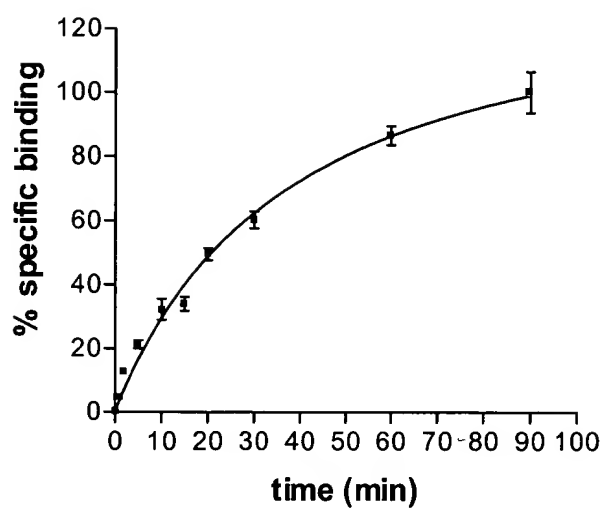


Figure 1: Association [3H]-4MG



101000-150000

Figure 2: Dissociation of  
[3H]-4MG binding

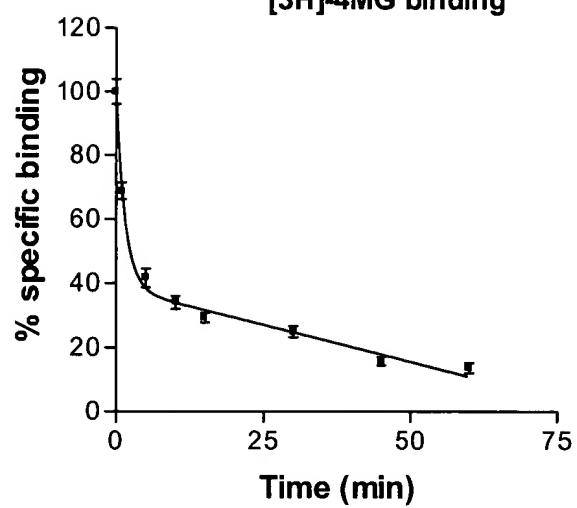


Figure 3: DRUG INHIBITION OF  
[3H]-4MG BINDING

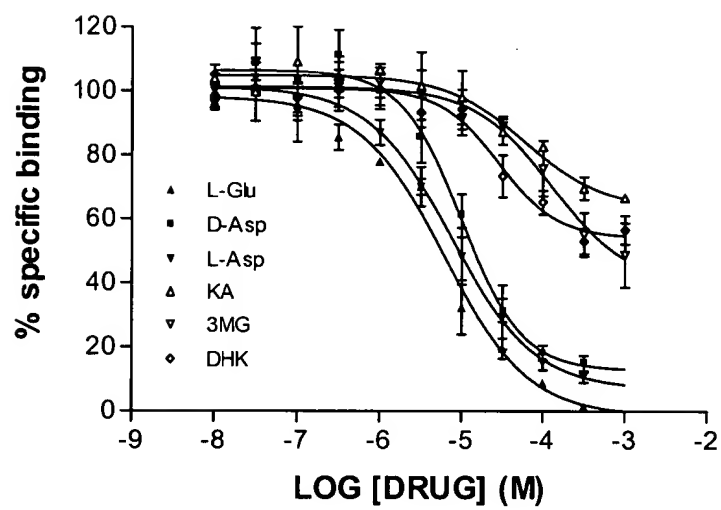


Figure 4: DRUG INHIBITION of [3H]-4MG BINDING

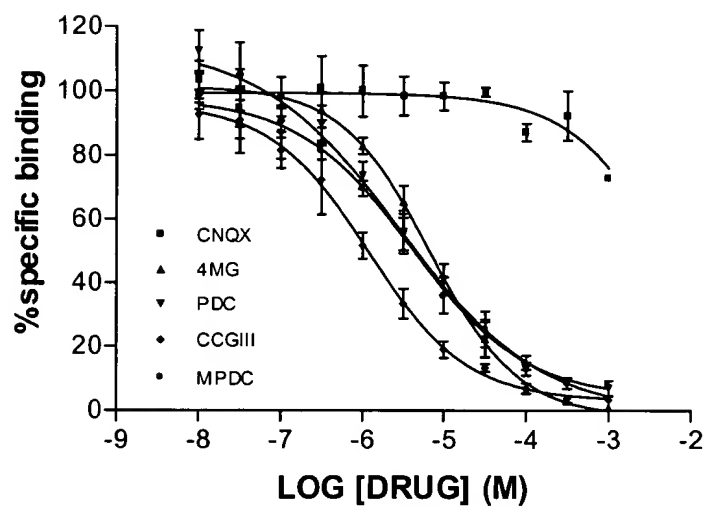
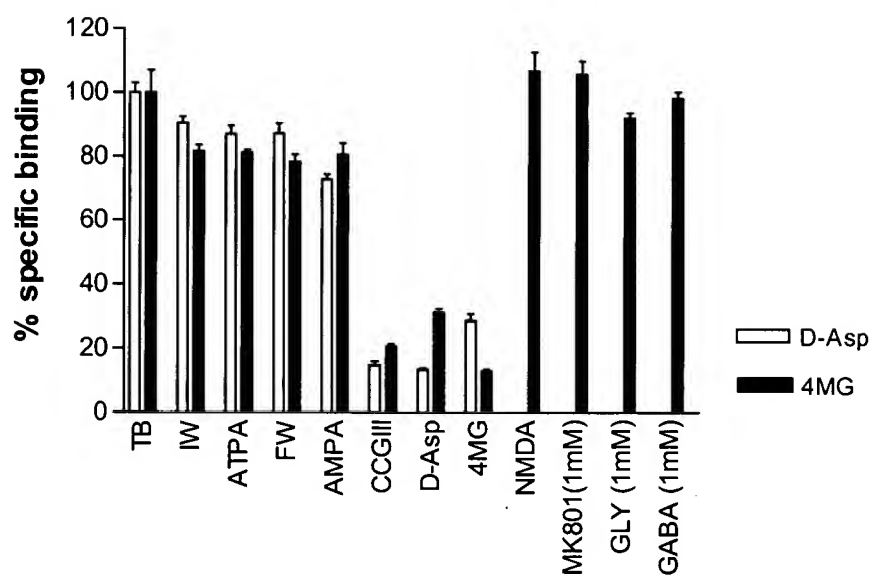


Figure 5: COMPARISON of MISCELLANEOUS DRUG  
INHIBITION of [3H]-D-ASPARTATE and [3H]-4MG



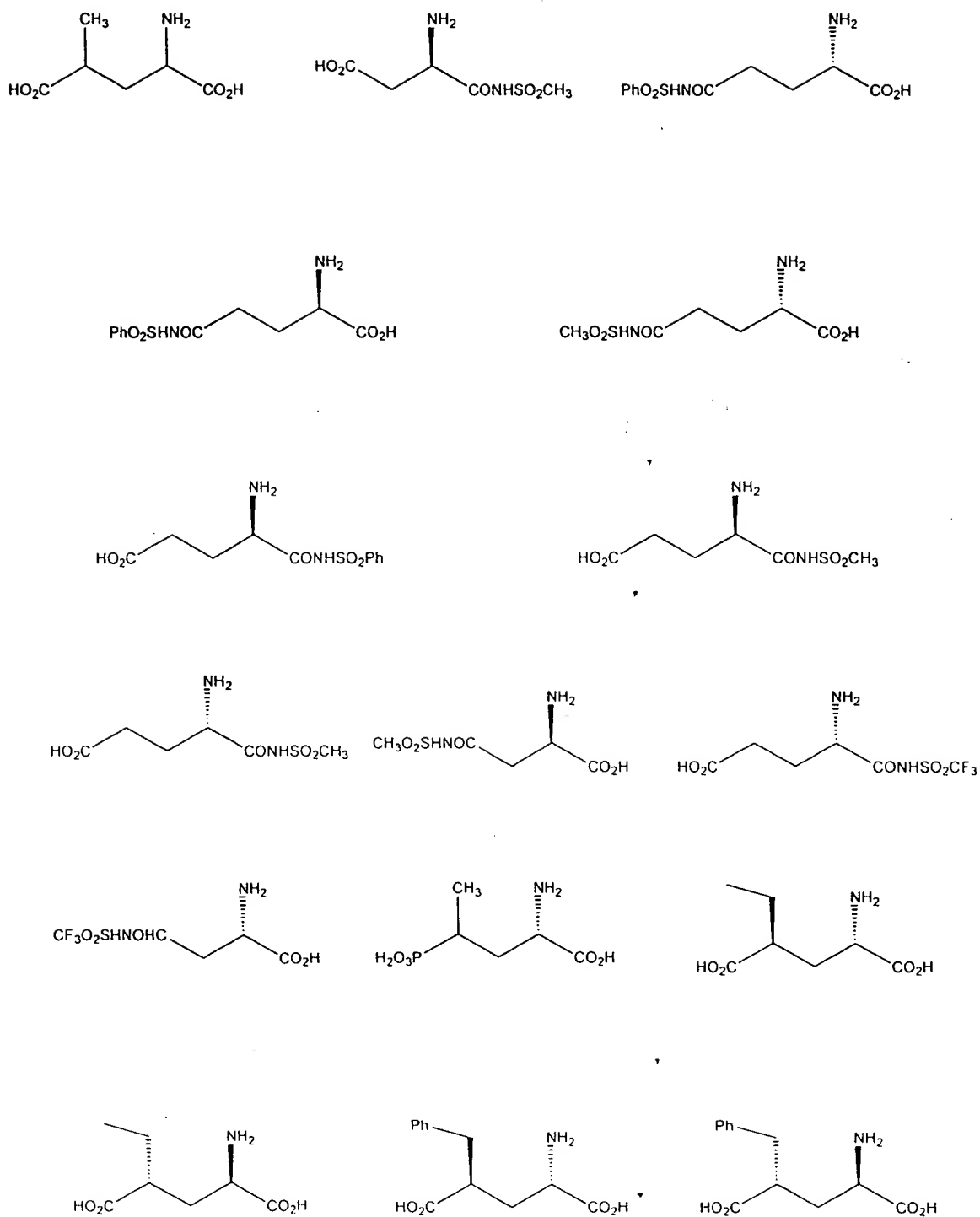


Figure 6A

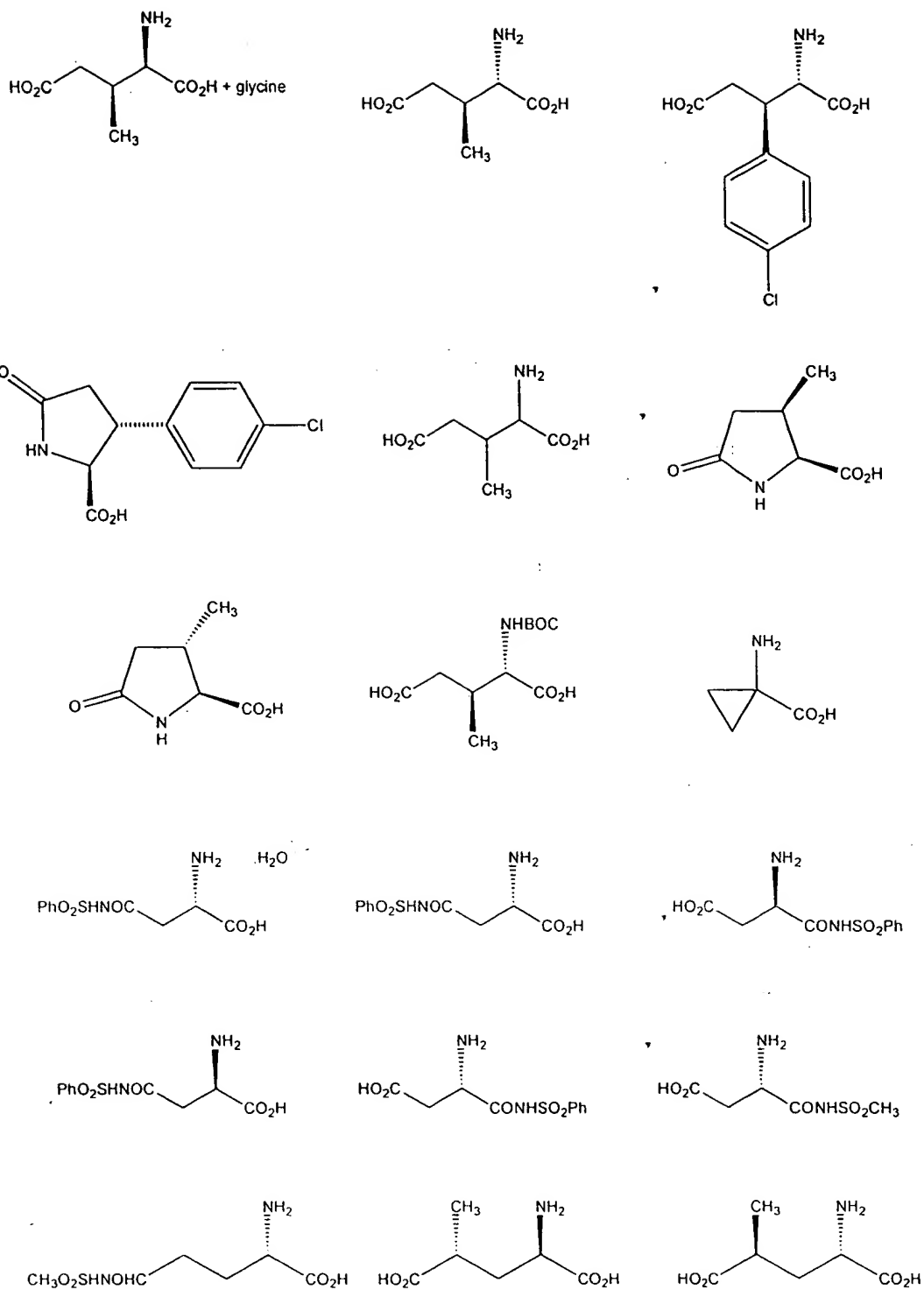


Figure 6B

OC(=O)[C@H](Cc1ccc(O)cc1)[C@@H](N)C(=O)O



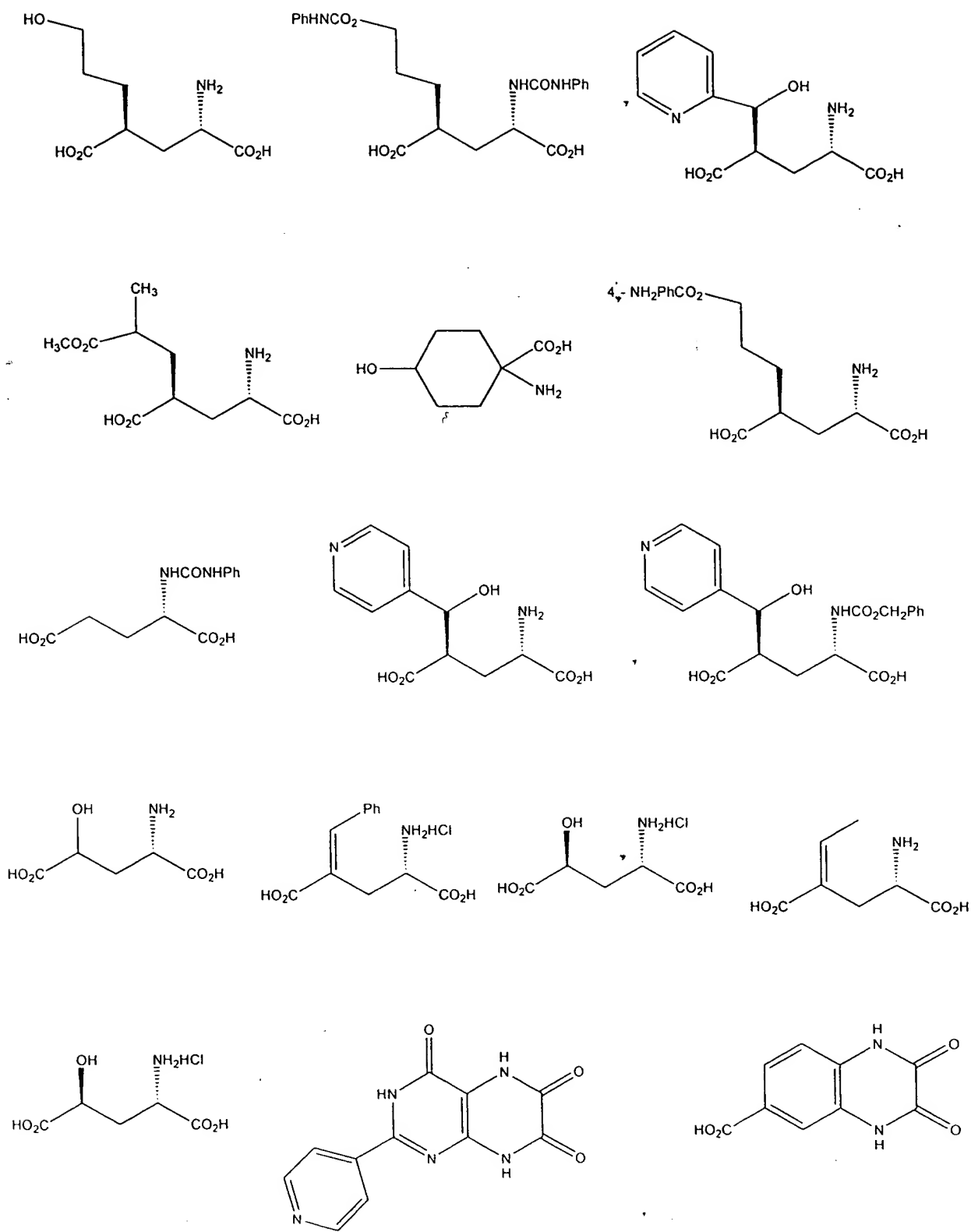


Figure 6D

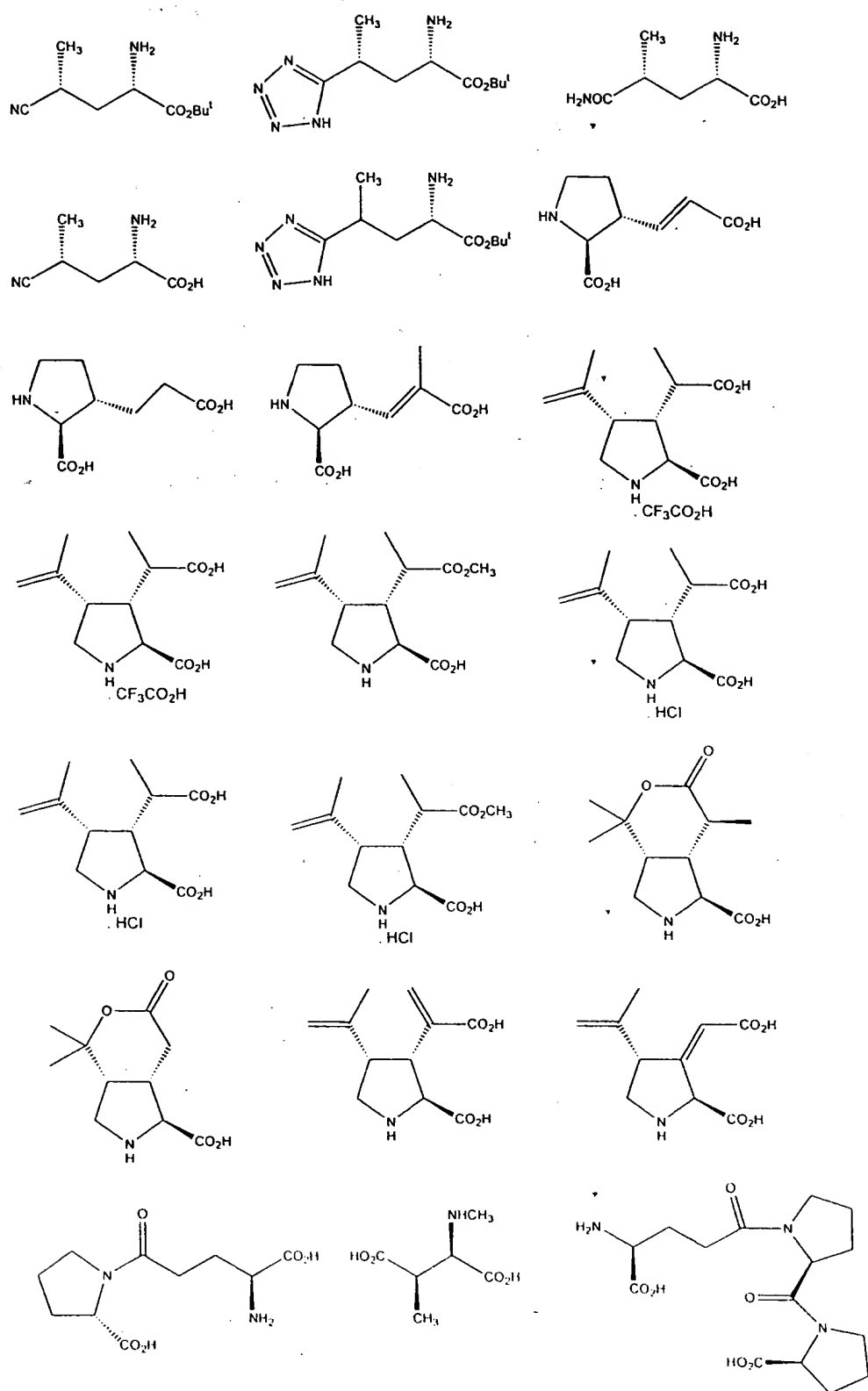


Figure 6E

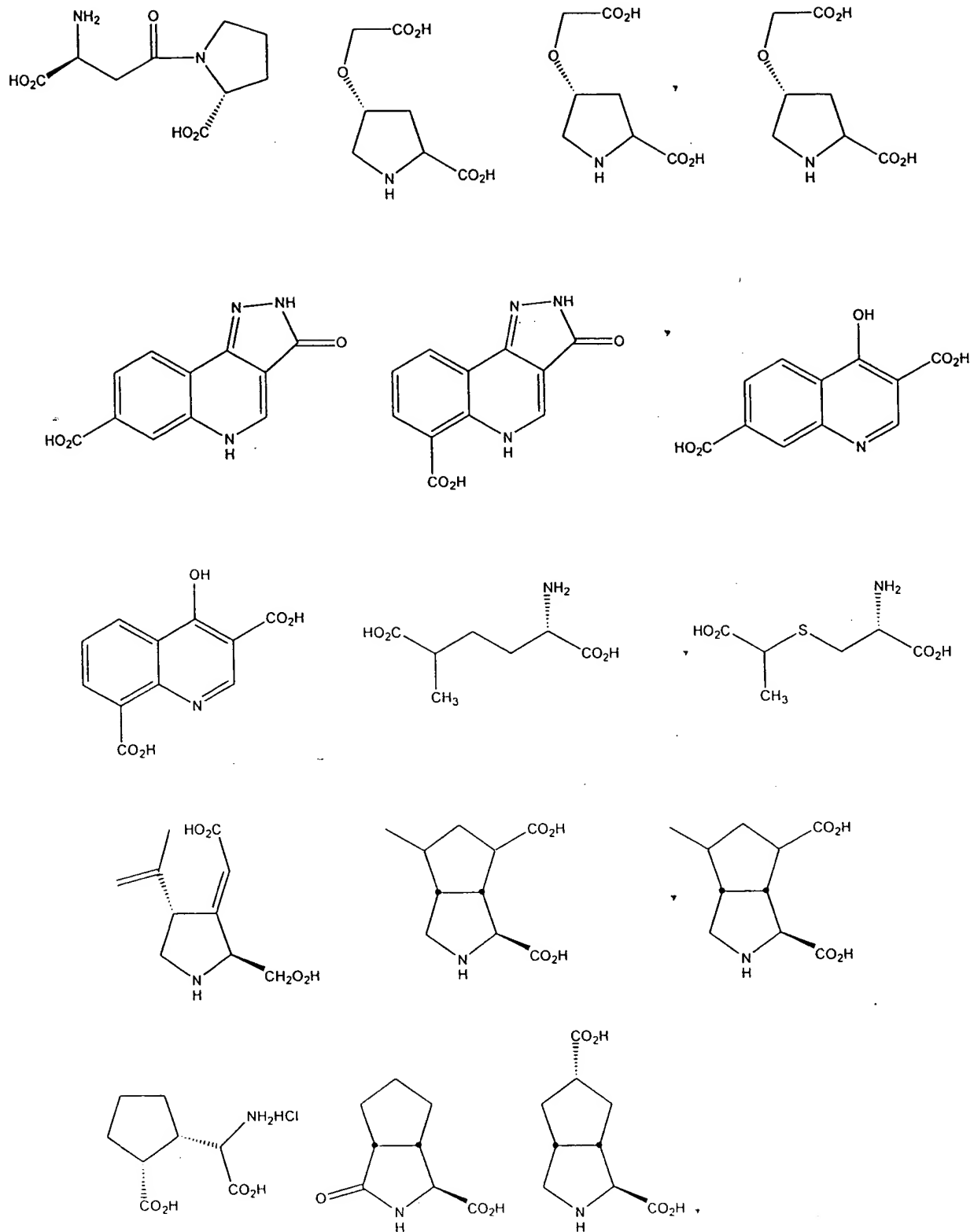


Figure 6F

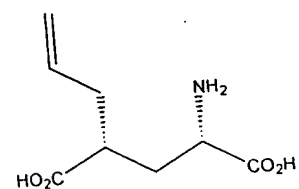
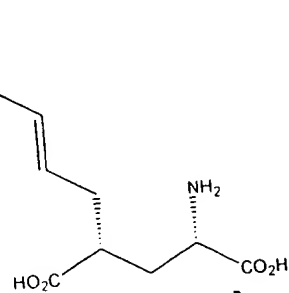
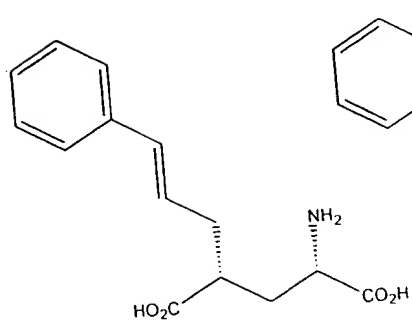
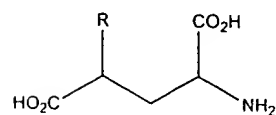
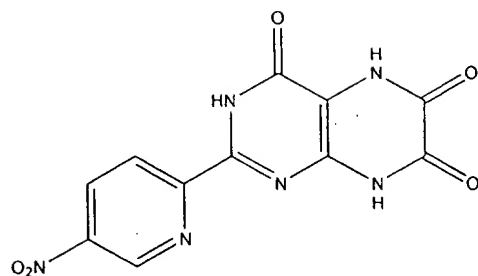
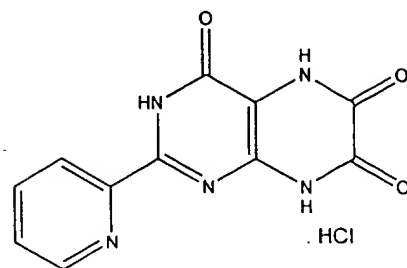
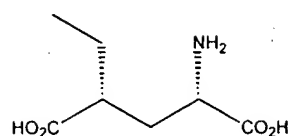
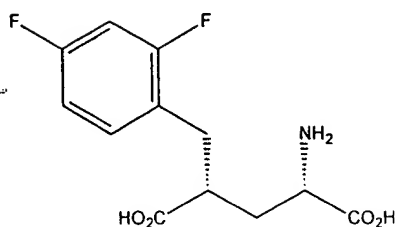
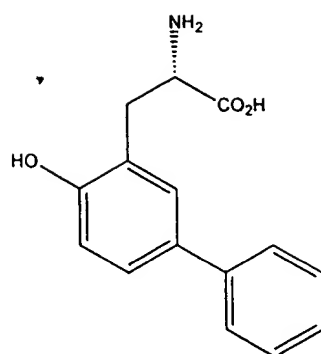
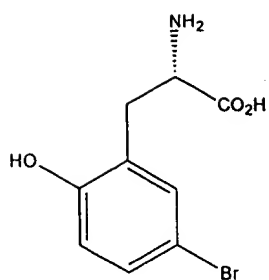
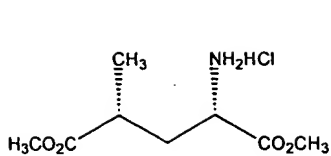


Figure 6G

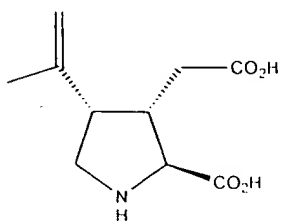
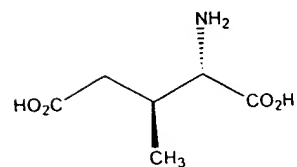
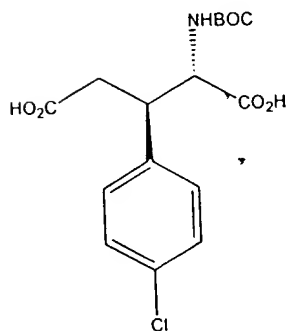
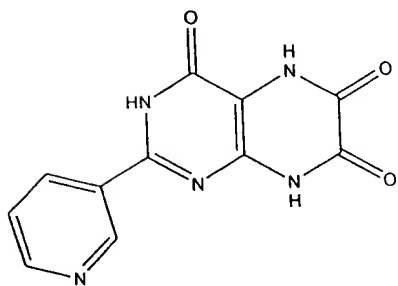
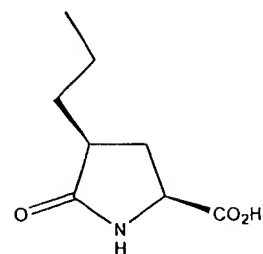
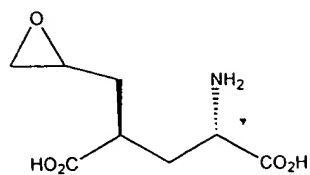
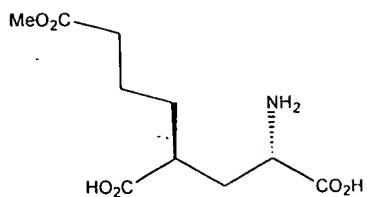
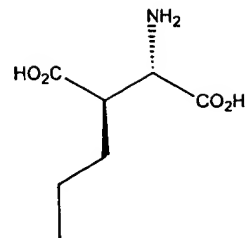
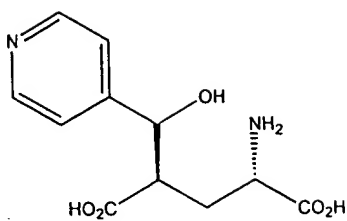
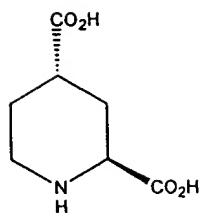
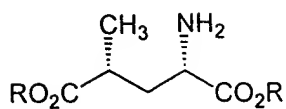
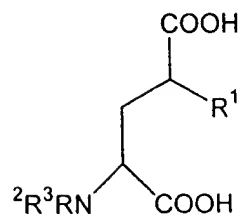
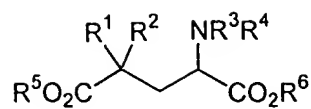


Figure 6H



R = Me, Et, <sup>t</sup>Bu

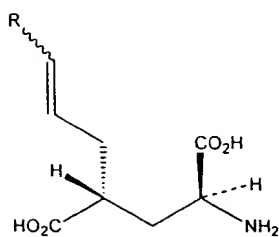
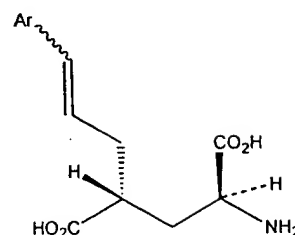


R<sup>1</sup> = CH<sub>3</sub>, and halogen

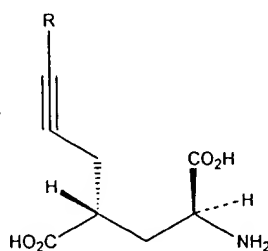
R<sup>2</sup>, R<sup>3</sup> are independently

H, C1-C6-alkyl, C3-C4-alkenyl, C3-C5-cycloalkyl, C1-C6-alkyl-CO-,  
C1-C6-alkyl-OCO-, C1-C6-alkyl-NHCO-, HCO-, or C3-C6-alkynyl

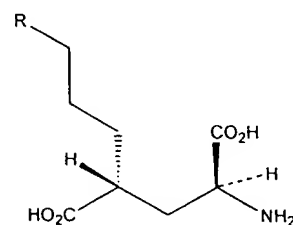
R<sup>2</sup>, R<sup>3</sup> taken together can be -CH<sub>2</sub>(CH<sub>2</sub>)<sub>p</sub>CH<sub>2</sub>-



R = H, Me, Et, Cl



R = H, Me, Et, nPr



R = H, Et, nPr

Figure 6I